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STRUCTURE FILE UPDATES: 25 MAR 2008 HIGHEST RN 1010115-69-1  
 DICTIONARY FILE UPDATES: 25 MAR 2008 HIGHEST RN 1010115-69-1

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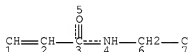
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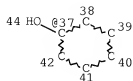
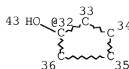
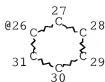
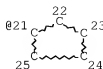
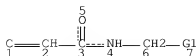
=> d sta que l74  
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 DEFAULT ECLEVEL IS LIMITED

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 NUMBER OF NODES IS 7

STEREO ATTRIBUTES: NONE  
 L60 42913 SEA FILE=REGISTRY SSS FUL L58  
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VAR G2=AK/ID/PH/13

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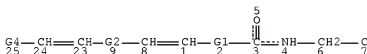
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L63 1639 SEA FILE=REGISTRY SUB=L60 CSS FUL L61

L72 STR



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VAR G4=AK/ID

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DEFAULT ECLEVEL IS LIMITED

GRAPH ATTRIBUTES:

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NUMBER OF NODES IS 14

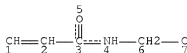
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L74 22 SEA FILE=REGISTRY SUB=L63 SSS FUL L72

100.0% PROCESSED 169 ITERATIONS  
 SEARCH TIME: 00.00.01

22 ANSWERS

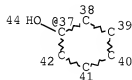
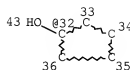
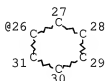
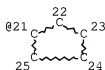
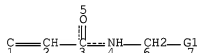
=> d sta que l66  
 L58 STR



NODE ATTRIBUTES:  
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 DEFAULT MLEVEL IS ATOM  
 DEFAULT ECLEVEL IS LIMITED

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 NUMBER OF NODES IS 7

STEREO ATTRIBUTES: NONE  
 L60 42913 SEA FILE=REGISTRY SSS FUL L58  
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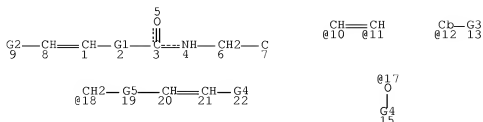
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 GGCAAT IS MCY UNS AT 14  
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 ECOUNT IS E6 C AT 14

GRAPH ATTRIBUTES:  
 RSPEC 21 26 32 37  
 NUMBER OF NODES IS 42

STEREO ATTRIBUTES: NONE

L63 1639 SEA FILE=REGISTRY SUB=L60 CSS FUL L61

L64 STR



REP G1=(0-1) 10-1 11-3

VAR G2=PH/12/18

VAR G3=AK/ID/17

VAR G4=AK/ID

REP G5=(0-2) CH2

NODE ATTRIBUTES:

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CONNECT IS M1 RC AT 7

DEFAULT MLEVEL IS ATOM

GGCAT IS MCY UNS AT 12

DEFAULT ECLEVEL IS LIMITED

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RING(S) ARE ISOLATED OR EMBEDDED

NUMBER OF NODES IS 20

STEREO ATTRIBUTES: NONE

L66 94 SEA FILE=REGISTRY SUB=L63 CSS FUL L64

100.0% PROCESSED 1639 ITERATIONS

94 ANSWERS

SEARCH TIME: 00.00.01

=&gt; fil hcaplus

FILE 'HCAPLUS' ENTERED AT 14:37:37 ON 26 MAR 2008

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FILE COVERS 1907 - 26 Mar 2008 VOL 148 ISS 13

FILE LAST UPDATED: 25 Mar 2008 (20080325/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

=> d bib abs hitind hitstr retrievable tot 1111

L111 ANSWER 1 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2005:983777 HCAPLUS Full-text

DN 143:266752

TI Processes for preparing (2E,4E,8Z)-2,4,8-undecatrienoic acid and ester and carboxamide derivatives and organoleptic uses thereof

IN Dewis, Mark L.; Huber, Michelle E.

PA International Flavors & Fragrances Inc., USA

SO U.S. Pat. Appl. Publ., 12 pp., Cont.-in-part of U.S. Ser. No. 618,367.

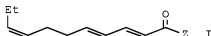
CODEN: USXXCO

DT Patent

LA English

FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2005197387	A1	20050908	US 2004-861751	20040604 <--
	US 7098350	B2	20060829		
	US 2005010062	A1	20050113	US 2003-618367	20030710 <--
	US 7141686	B2	20061128		
	IN 2004DE01233	A	20060721	IN 2004-DE1233	20040701 <--
	EP 1496042	A2	20050112	EP 2004-254095	20040708 <--
	EP 1496042	A3	20050309		
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK, HR				
	CN 1706788	A	20051214	CN 2004-10063646	20040709 <--
PRAI	US 2003-618367	A2	20030710	<--	
	US 2004-861751	A	20040604		
OS	CASREACT 143:266752; MARPAT 143:266752				
GI					

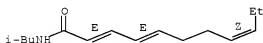


AB Described is a genus of undecatrienoic acid derivs. I [Z = NRR1, OR2; R = H, Me, Et; R1 = Me, Et, Pr, cyclopropyl, iso-Pr, Bu, s-Bu, iso-Bu, 2-methylbutyl, cyclobutyl, piperonyl, cyclophenyl, allyl; R2 = H, (un)branched C1-6-alkyl, C3-6-alkenyl] useful in imparting, augmenting and/or enhancing flavors, aromas and somatosensory effects in or to consumable materials such as foods, beverages, skin care products, oral care products, medicinal products and the like. Also described is a synthesis process for producing such derivs. The process comprises: (a) oxidation of (2E,4E,8Z)-2,4,8-undecatrienal with Ag2O in aqueous alkali hydroxide, followed by acidifying the product; (b) reacting the acid with an alkyl haloformate in the presence of a tertiary amine; (c) reacting the intermediate mixed anhydride with either an amine, RR1NH, to form the amide, or reacting with an alc., R2OH, to form the ester. Thus, (2E,4E,8Z)-N-(isobutyl)-2,4,8-undecatrienamide [I; Z = NHCH2CHMe2], was prepared from (2E,4E,8Z)-2,4,8-undecatrienal [Z = H] via oxidation with Ag2O in aqueous NaOH, acidification with aqueous HCl, reaction with ClCO3Et in the presence of Et3N, then amination with Me2CHCH2NH2.

IC ICM A61K0031-202

- ICS A61K0031-16; A61K0031-20  
 INCL 514464000; X51-456.0; X51-462.7; X55-4 3.5; X55-422.3  
 CC 26-3 (Biomolecules and Their Synthetic Analogs)  
 Section cross-reference(s): 17, 62, 63  
 IT Beverages  
 Candy  
 Chewing gum  
 Drugs  
 (flavor enhancers for; preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 IT Condiments  
 (flavor-enhancing; preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 IT Dentistry  
 (oral care products, flavor enhancers for; preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 IT Food additives  
 Human  
 Odor and Odorous substances  
 Taste  
 (preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 IT Cooperative phenomena  
 (synergism, with aroma, taste or somatosensory agents; preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 IT 652970-05-3P, (2E,4E,8Z)-N-(Isobutyl)-2,4,8-undecatrienamide  
 RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
 (preparation and organoleptic uses of; preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 IT 652970-05-3P, (2E,4E,8Z)-N-(Isobutyl)-2,4,8-undecatrienamide  
 RL: PRP (Properties); SPN (Synthetic preparation); PREP (Preparation)  
 (preparation and organoleptic uses of; preparation of (2E,4E,8Z)-2,4,8-undecatrienoic acid ester and carboxamide derivs. and their organoleptic uses)  
 RN 652970-05-3 HCAPLUS  
 CN 2,4,8-Undecatrienamide, N-(2-methylpropyl)-, (2E,4E,8Z)- (CA INDEX NAME)

Double bond geometry as shown.



## RETABLE

Referenced Author (RAU)	Year (RPY)	VOL (RVL)	PG (RPG)	Referenced Work (RWK)	Referenced File
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Anon	1970			JP 04803546	
Anon	1976			GB 1438205	HCAPLUS
Anon	1981			JP 56087505	HCAPLUS
Anon	1993			WO 9323005	HCAPLUS
Anon	1998			WO 9807404	HCAPLUS
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Anon	2000			W0 0/45815	HCAPLUS
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Anon	2002			W0 0/051392	
Anon	2003			W0 2004000787 A2	HCAPLUS
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Anon	2001	55	53	Food Technology	
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Anon				Search for Unsaturat	
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Gamboa-Leon, R	2000	28	1019	Biochemical Systemat	HCAPLUS
Gamboa-Leon, R	2000	28	1019	Biochemical Systemat	HCAPLUS
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Nakatsu	1996			US 5545424 A	HCAPLUS
Ottinger	2001	49	5383	J. Agric. Food Chem.	HCAPLUS
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Saadali, B	2001  58	1083	Phytochemistry	HCAPLUS
Sako	1999		US 5955066 A	HCAPLUS
Shiroyama	2001		US 6328982 B1	HCAPLUS
Sonnenberg	2002		US 20020173436 A1	HCAPLUS
Tashjian	2003		US 6579513 B1	HCAPLUS
Valentine	2003		US 6579535 B1	HCAPLUS
Vermeer	1997		US 5641480 A	HCAPLUS
Watkins	2002		US 6451844 B1	HCAPLUS
Watson	1979		US 4150052 A	HCAPLUS
Watson	1980		US 4226988 A	HCAPLUS
Winkler	1984		US 4470982 A	HCAPLUS
Wolf	2002		US 6455080 B1	HCAPLUS
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Yeoh	2001		US 6200554 B1	HCAPLUS
Young	2001		US 6248315 B1	HCAPLUS

L111 ANSWER 2 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2005:34644 HCAPLUS [Full-text](#)

DN 142:134213

TI Ester and carboxamide derivatives of E2,E4,Z8-undecatienoic acid, processes for preparing same and their organoleptic uses

IN Lewis, Mark L.; Huber, Michelle E.

PA International Flavors & Fragrances Inc., USA

SO U.S. Pat. Appl. Publ., 10 pp.

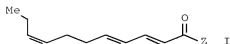
CODEN: USXXCO

DT Parent

LA English

FAN.CNT 2

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	US 7141686	B2	20061128		
	US 2005197387	A1	20050908	US 2004-861751	20040604 <--
	US 7098350	B2	20060829		
	IN 2004DE01233	A	20060721	IN 2004-DE1233	20040701 <--
	EP 1496042	A2	20050112	EP 2004-254095	20040708 <--
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	CN 1706788	A	20051214	CN 2004-10063646	20040709 <--
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GI					

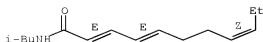




- AB The present invention discloses preparation of E2,E4,Z8-undecatrienoic acid derivs., such as I [Z = OR1, NR2R3; R1 = H, C1-C6 straight chain or branched-chain alkyl, C3-C6 straight chain or branched-chain alkenyl; R2 = H, Me, Et; R3 = Me, Et, Pr, cyclopropyl, iso-Pr, Bu, sec-Bu, iso-Bu, 2-methylbutyl, cyclobutyl, 3,4-methylenedioxyphenyl, cyclopentyl or allyl], and their use in imparting, augmenting and/or enhancing flavors, aromas and somatosensory effects in or to consumable materials such as foods, beverages, skin care products, oral care products, medicinal products and the like. Thus, I [Z = NHCH2CHMe2 (II)] was prepared by the reaction of isobutylamine and E2,E4,Z8-undecatrienoic acid (prepared via oxidation of E2,E4,Z8-undecatrilal with silver(I) oxide). II was used for the enhancement of flavor of alc. beverage and hard candy.
- IC ICM C07C0233-02  
ICS C07D0317-44
- INCL 554035000; X55-422.4; X54-943.6
- CC 23-18 (Aliphatic Compounds)  
Section cross-reference(s): 17, 62, 63
- ST undecatrienoic acid ester carboxamide deriv prepn food cosmetic additive; beverage candy chewing gum toothpaste additive undecatrienoic acid deriv
- IT Carboxylic acids, preparation  
RL: COS (Cosmetic use); FFD (Food or feed use); IMF (Industrial manufacture); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(derivs.; preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT Carboxylic acids, preparation  
RL: COS (Cosmetic use); FFD (Food or feed use); IMF (Industrial manufacture); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(esters; preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT Candy  
(hard; preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT Drug delivery systems  
(nasal; preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT Hygiene  
(oral; preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT Alcoholic beverages  
Chewing gum  
Colognes  
Cosmetics  
Dentifrices  
Flavoring materials  
Hair preparations  
Skin preparations (pharmaceutical)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT Amides, preparation

- RL: COS (Cosmetic use); FFD (Food or feed use); IMF (Industrial manufacture); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 652970-05-3P  
RL: COS (Cosmetic use); FFD (Food or feed use); IMF (Industrial manufacture); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 824416-99-1P, E2,E4,Z8-Undecatrienoic acid  
RL: IMF (Industrial manufacture); RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 795309-53-4P 823815-34-5P  
RL: IMF (Industrial manufacture); SPN (Synthetic preparation); PREP (Preparation)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 67-56-1, Methanol, reactions 78-81-9, Iso-butylamine 124-40-3, Dimethylamine, reactions 541-41-3, Ethyl chloroformate 350696-20-7, E2,E4,Z8-Undecatrienal  
RL: RCT (Reactant); RACT (Reactant or reagent)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 823815-35-6P  
RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 121-44-8, Triethylamine, reactions 1310-73-2, Sodium hydroxide, reactions 7647-01-0, Hydrochloric acid, reactions 20667-12-3, Silver(I) oxide  
RL: RGT (Reagent); RACT (Reactant or reagent)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- IT 652970-05-3P  
RL: COS (Cosmetic use); FFD (Food or feed use); IMF (Industrial manufacture); SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation of ester and carboxamide derivs. of E2,E4,Z8-undecatrienoic acid and their uses for enhancing flavors, aromas and somatosensory effects in consumable materials)
- RN 652970-05-3 HCAPLUS  
CN 2,4,8-Undecatrienamide, N-(2-methylpropyl)-, (2E,4E,8Z)- (CA INDEX NAME)

Double bond geometry as shown.



## RETABLE

Referenced Author (RAU)	Year (RPY)	VOL (RVL)	PG (RPG)	Referenced Work (RWK)	Referenced File
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Anon	1981			JP 56087505	HCAPLUS
Anon	1993			WO 9323005	HCAPLUS
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Anon	2004			WO 2004011415	HCAPLUS
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Gamboa-Leon	2000	28	1019	Biochemical Systemat	HCAPLUS
Gatfield	2004			US 20040241312 A1	
Glenn	2003			US 6544499 B1	HCAPLUS
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Grainger	2002			US 6365215 B1	HCAPLUS
Guskey	2001			US 6297203 B1	HCAPLUS
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Hammer	2003			US 20030082124 A1	HCAPLUS
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Jacobson	1953	75	2584	Pellitorine Isomers.	HCAPLUS
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Jarboe	1963			US 3111127 A	HCAPLUS
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Rapaport	1998			US 5730965 A	HCAPLUS
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Rowseell	1977			US 4032661 A	HCAPLUS
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Tashjian	2003			US 6579513 B1	HCAPLUS
Valentine	2003			US 6579535 B1	HCAPLUS
Vermeer	1997			US 5641480 A	HCAPLUS
Watkins	2002			US 6451844 B1	HCAPLUS
Watson	1979			US 4150052 A	HCAPLUS
Watson	1980			US 4226988 A	HCAPLUS
Winkler	1984			US 4470982 A	HCAPLUS
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Wolf	2003			US 20030082271 A1	
Wolfson	2002			US 20020122778 A1	HCAPLUS
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Young	2001			US 6248315 B1	HCAPLUS
Zimmermann	2001			US PP12213 P2	

L111 ANSWER 3 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2094:673840 HCAPLUS [Full-text](#)

DN 141:325774

TI Sanshool derivatives from Zanthoxylum piperitum as memory enhancers and health foods

IN Yano, Shingo; Nakamura, Tomonori; Ikegami, Fumio

PA Tokiwa Shikubutsu Kagaku Kenkyusho Ltd., Japan

SO Jpn. Kokai Tokkyo Koho, 17 pp.

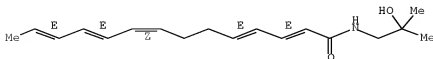
CODEN: JKXXAF

DT Patent

LA Japanese  
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 2004292383	A	20041021	JP 2003-88262	20030327 <--
PRAI	JP 2003-88262		20030327	<--	
OS	MARPAT 141:325774				
AB	Zanthol derivs. from Zanthoxylum piperitum (I; Me(CH <sub>2</sub> ) <sub>n</sub> CONHCH <sub>2</sub> R <sub>1</sub> wherein R <sub>1</sub> = Me, CH <sub>2</sub> OR <sub>2</sub> , CR <sub>3</sub> (CH <sub>3</sub> ) <sub>2</sub> , etc., with R <sub>2</sub> =, H, Me, sugar, R <sub>3</sub> = H, OH) are claimed as memory enhancers and health foods. I were extracted from the above plant, and their effects on learning were studied in mouse water maze test.				
IC	ICM A61K0031-16 ICS A23L0001-30; A61K0031-164; A61K0035-78; A61P0025-28				
CC	1-11 (Pharmacology) Section cross-reference(s): 17				
ST	sanshool deriv Zanthoxylum memory enhancer health food				
IT	Cognition enhancers Health food Learning Zanthoxylum piperitum (sanshool derivs. from Zanthoxylum piperitum as memory enhancers and health foods)				
IT	10076-00-3P, β-Sanshool 78886-65-4P, γ-Sanshool 78886-66-5E, Hydroxy-γ-Sanshool 97465-69-5E, Hydroxy-β-Sanshool RL: PAC (Pharmacological activity); PUR (Purification or recovery); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (sanshool derivs. from Zanthoxylum piperitum as memory enhancers and health foods)				
IT	504-97-2, α-Sanshool 83883-10-7, Hydroxy-α-Sanshool RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL (Biological study); USES (Uses) (sanshool derivs. from Zanthoxylum piperitum as memory enhancers and health foods)				
IT	78886-66-5P, Hydroxy-γ-Sanshool 97465-69-5P, Hydroxy-β-Sanshool RL: PAC (Pharmacological activity); PUR (Purification or recovery); THU (Therapeutic use); BIOL (Biological study); PREP (Preparation); USES (Uses) (sanshool derivs. from Zanthoxylum piperitum as memory enhancers and health foods)				
RN	78886-66-5 HCAPLUS				
CN	2,4,8,10,12-Tetradecapentaenamide, N-(2-hydroxy-2-methylpropyl)-, (2E,4E,8Z,10E,12E)- (CA INDEX NAME)				

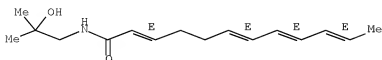
Double bond geometry as shown.



RN 97465-69-5 HCAPLUS

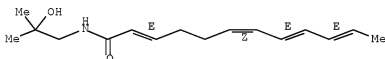
CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-, (2E,6E,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



IT 93883-10-7, Hydroxy- $\alpha$ -Sanshool  
 RL: PAC (Pharmacological activity); THU (Therapeutic use); BIOL  
 (Biological study); USES (Uses)  
 (sanshool derivs. from Zanthoxylum piperitum as memory enhancers and  
 health foods)  
 RN 93883-10-7 HCAPLUS  
 CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-,  
 (2E,6Z,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



L111 ANSWER 4 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2004:856961 HCAPLUS [Full-text](#)

DN 141:331119

TI Alkylidienamides exhibiting taste and sensory effect in  
 flavor compositions

IN Dewis, Mark L.; Huber, Michelle E.; Cossette, Michael V.; Agyemang, David  
 O.

PA International Flavors & Fragrances Inc, USA

SO U.S. Pat. Appl. Publ., 9 pp.

CODEN: USXXCO

DT Patent

LA English

FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2004202760	A1	20041014	US 2003-411672	20030411 <--
	US 2004202619	A1	20041014	US 2004-783652	20040220 <--
	IN 2004DE00651	A	20060616	IN 2004-DE651	20040331 <--
	BR 2004001566	A	20050830	BR 2004-1566	20040406 <--
	EP 1473287	A2	20041103	EP 2004-252136	20040408 <--
	EP 1473287	A3	20041229		
	EP 1473287	B1	20060621		
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK, HR				
	ES 2267004	T3	20070301	ES 2004-252136	20040408 <--
	CN 1593230	A	20050316	CN 2004-10032498	20040409 <--
PRAI	US 2003-411672	A2	20030411	<--	
	US 2004-783652	A	20040220		
OS	MARPAT 141:331119				

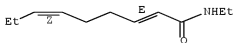
- AB Alkyldienamide compds. suitable for use as flavoring agents are disclosed. The compds. are used as flavors since they possess umami characteristics or other desirable organoleptic properties.
- IC ICM A23L0001-22
- INCL 426534000
- CC 17-6 (Food and Feed Chemistry)  
Section cross-reference(s): 23, 24, 62
- ST alkyldienamide flavor enhancer synthesis food oral  
hygiene additive; nonadienamide dodecadienamide deriv flavor  
enhancer food toothpaste
- IT Beverages  
Chewing gum  
Dentifrices  
Flavoring materials  
Food additives  
(alkyldienamides exhibiting taste and sensory effect in  
flavor compns.)
- IT Amides, biological studies  
RL: FFD (Food or feed use); BIOL (Biological study); USES (Uses)  
(alkyldienes; alkyldienamides exhibiting taste and sensory  
effect in flavor compns.)
- IT Dentifrices  
(aromatic oil-flavored; alkyldienamides exhibiting taste  
and sensory effect in flavor compns.)
- IT Chewing gum  
(bubble gum flavored; alkyldienamides exhibiting  
taste and sensory effect in flavor compns.)
- IT Condiments  
(flavor-enhancing; alkyldienamides exhibiting taste  
and sensory effect in flavor compns.)
- IT Candy  
(hard, cinnamon-flavored; alkyldienamides exhibiting  
taste and sensory effect in flavor compns.)
- IT Beverages  
(lemon-lime flavor; alkyldienamides exhibiting taste  
and sensory effect in flavor compns.)
- IT Hygiene  
(oral, products for; alkyldienamides exhibiting taste and  
sensory effect in flavor compns.)
- IT Alcoholic beverages  
(peach flavor; alkyldienamides exhibiting taste and  
sensory effect in flavor compns.)
- IT 608514-55-2P 608514-56-3P 767329-59-9P 767329-60-2P  
767329-61-3P 767329-62-4P 767329-65-7P 767329-66-8P  
767329-68-0P 767329-69-1P 767329-70-4P 767329-71-5P 767329-72-6P  
773060-63-2P 773060-64-3DP, derivs. 773060-65-4DP, derivs.  
RL: FFD (Food or feed use); SPN (Synthetic preparation); BIOL  
(Biological study); PREP (Preparation); USES (Uses)  
(alkyldienamides exhibiting taste and sensory effect in  
flavor compns.)
- IT 74-89-5, Methylamine, reactions 75-04-7, Ethylamine, reactions  
75-31-0, Isopropylamine, reactions 78-81-9, Isobutylamine 96-15-1,  
2-Methylbutylamine 141-43-5, 2-Ethanolamine, reactions 541-41-3, Ethyl  
chloroformate 765-30-0, Cyclopropylamine 2620-50-0, Piperonylamine  
23605-13-2 94088-26-3  
RL: RCT (Reactant); RACT (Reactant or reagent)  
(alkyldienamides exhibiting taste and sensory effect in  
flavor compns.)
- IT 608514-56-3P 767329-59-9P 767329-61-3P  
773060-63-2P

RL: EFD (Food or feed use); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses) (alkyldienamides exhibiting taste and sensory effect in flavor compns.)

RN 608514-56-3 HCAPLUS

CN 2,6-Nonadienamide, N-ethyl-, (2E,6Z)- (CA INDEX NAME)

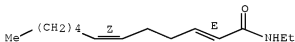
Double bond geometry as shown.



RN 767329-59-9 HCAPLUS

CN 2,6-Dodecadienamide, N-ethyl-, (2E,6Z)- (CA INDEX NAME)

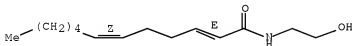
Double bond geometry as shown.



RN 767329-61-3 HCAPLUS

CN 2,6-Dodecadienamide, N-(2-hydroxyethyl)-, (2E,6Z)- (CA INDEX NAME)

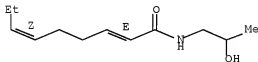
Double bond geometry as shown.



RN 773060-63-2 HCAPLUS

CN 2,6-Nonadienamide, N-(2-hydroxypropyl)-, (2E,6Z)- (CA INDEX NAME)

Double bond geometry as shown.



L111 ANSWER 5 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2004:856915 HCAPLUS [Full-text](#)

DN 141:313274

TI Compositions comprising alkyldienamides exhibiting taste and sensory effect and use for enhancing flavor in foodstuff and beverage

IN Lewis, Mark L.; Huber, Michelle E.; Cossette, Michael V.; Agyemang, David O.; Conklin, Garry; Pei, Tao

PA International Flavors & Fragrances Inc, USA



SO U.S. Pat. Appl. Publ., 15 pp., Cont.-in-part of Ser. No. US 2003-411672,  
filed on 11 Apr 2003  
CODEN: USXXCO

DT Patent

LA English

FAN.CNT 2

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2004202619	A1	20041014	US 2004-783652	20040220 <--
	US 2004202760	A1	20041014	US 2003-411672	20030411 <--
	BR 2004001566	A	20050830	BR 2004-1566	20040406 <--
	EP 1473287	A2	20041103	EP 2004-252136	20040408 <--
	EP 1473287	A3	20041229		
	EP 1473287	B1	20060621		
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK, HR				
	ES 2267004	T3	20070301	ES 2004-252136	20040408 <--
	CN 1593230	A	20050316	CN 2004-10032498	20040409 <--
PRAI	US 2003-411672	A2	20030411	<--	
	US 2004-783652	A	20040220		

OS MARPAT 141:313274

AB The present invention relates to novel compds. and a process for augmenting or imparting a taste or somatosensory effect to a foodstuff, chewing gum, medicinal product, toothpaste, alc. beverage, aqueous beverage or soup. The compds. are used as flavors since they possess umami characteristics or other desirable organoleptic properties. The disclosed compds. are defined by the formula R5R4C:C(R3)CON(Y)(X) (X = H, Me, Et, Pr, iPr; Y = Me, Et, cyclopropyl, iPr, Pr, Bu, sec-Bu, iso-Bu, 2-methylbutyl, allyl, cyclobutyl, cyclopentyl, CH2CH(OH)CH3, CH(CH3)CH2OH, CH2C(CH3)OH, CH2CH2OH, CH2CO2CH3, geranyl, neryl, benzol[1,3]dioxol-5-yl; or X and Y together form the structures pyrrolidin-1-yl, 2-carboxypyrrolidin-1-yl, piperidin-1-yl; R3, R4 = Me, H; R5 = Me, Ph, benzyl, Et, Pr, Bu, iPr, phenylethyl, etc.).

IC ICM A61K0009-68

ICS A61K0007-16; A23L0001-221

INCL 424048000; X42-4 4.9; X42-665.0

CC 17-6 (Food and Feed Chemistry)

ST alkylidenamide taste sensory effect enhancing flavor  
foodstuff beverage

IT Alcoholic beverages

Chewing gum

Dentifrices

Flavor

Flavoring materials

Food

Soups

(alkylidenamides exhibiting taste and sensory effect and use for enhancing flavor in foodstuff and beverage)

IT Beverages

(aqueous; alkylidenamides exhibiting taste and sensory effect and use for enhancing flavor in foodstuff and beverage)

IT Candy

(hard; alkylidenamides exhibiting taste and sensory effect and use for enhancing flavor in foodstuff and beverage)

IT Saltiness

(salt taste enhancer; alkylidenamides exhibiting taste and sensory effect and use for enhancing flavor in foodstuff and beverage)

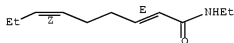
IT Taste

(somatosensory effect; alkylidenamides exhibiting taste and

sensory effect and use for enhancing flavor in  
foodstuff and beverage)

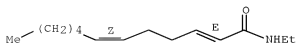
- IT 142-47-2, Monosodium glutamate 7647-14-5, Salt, biological studies  
80702-47-2, Ribotide  
RL: FFD (Food or feed use); BIOL (Biological study); USES (Uses)  
(alkyldienamides exhibiting taste and sensory effect and use  
for enhancing flavor in foodstuff and beverage)
- IT 608514-55-2P 608514-56-3P 767329-58-8P 767329-59-9P  
767329-60-2P 767329-61-3P 767329-62-4P 767329-63-5P  
767329-64-6P 767329-65-7P 767329-66-8P 767329-67-9P  
RL: FFD (Food or feed use); SPN (Synthetic preparation); BIOL  
(Biological study); PREP (Preparation); USES (Uses)  
(alkyldienamides exhibiting taste and sensory effect and use  
for enhancing flavor in foodstuff and beverage)
- IT 56-40-6, Glycine, reactions 74-89-5, Methylamine, reactions 75-04-7,  
Ethylamine, reactions 75-31-0, Isopropylamine, reactions 78-81-9,  
Isobutylamine 96-15-1, 2-Methylbutylamine 107-11-9, Allylamine  
121-44-8, Triethylamine, reactions 124-40-3, Dimethylamine, reactions  
141-43-5, 2-Ethanolamine, reactions 541-41-3, Ethyl chloroformate  
541-47-9 765-30-0, Cyclopropylamine 2620-50-0, Piperonylamine  
4698-08-2 23605-13-2 55320-96-2 75091-79-1 94088-26-3  
767329-75-9  
RL: RCT (Reactant); RACT (Reactant or reagent)  
(alkyldienamides exhibiting taste and sensory effect and use  
for enhancing flavor in foodstuff and beverage)
- IT 51552-27-3P 767329-68-0P 767329-69-1P 767329-70-4P 767329-71-5P  
767329-72-6P 767329-73-7P 767329-74-8P 767329-76-0P 767329-77-1P  
767329-78-2P 767329-79-3P 767329-80-6P 767329-81-7P  
RL: SPN (Synthetic preparation); PREP (Preparation)  
(alkyldienamides exhibiting taste and sensory effect and use  
for enhancing flavor in foodstuff and beverage)
- IT 608514-56-3P 767329-59-9P 767329-61-3P  
RL: FFD (Food or feed use); SPN (Synthetic preparation); BIOL  
(Biological study); PREP (Preparation); USES (Uses)  
(alkyldienamides exhibiting taste and sensory effect and use  
for enhancing flavor in foodstuff and beverage)
- RN 608514-56-3 HCAPLUS  
CN 2,6-Nonadienamide, N-ethyl-, (2E,6Z)- (CA INDEX NAME)

Double bond geometry as shown.



- RN 767329-59-9 HCAPLUS  
CN 2,6-Dodecadienamide, N-ethyl-, (2E,6Z)- (CA INDEX NAME)

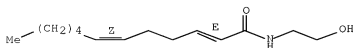
Double bond geometry as shown.



- RN 767329-61-3 HCAPLUS

CN 2,6-Dodecadienamide, N-(2-hydroxyethyl)-, (2E,6Z)- (CA INDEX NAME)

Double bond geometry as shown.



L111 ANSWER 6 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2004:428897 HCAPLUS Full-text

DN 141:6843

TI Preparation and use of trans-pellitorin as an aromatic substance with salivation-stimulating activity.

IN Gatfield, Ian Lucas; Ley, Jakob Peter; Krammer, Gerhard; Bertram, Heinz-Juergen; Loenneker, Ilse; Machinek, Arnold

PA Symrise G.m.b.H. &amp; Co. K.-G., Germany

SO PCT Int. Appl., 22 pp.

CODEN: PIXXD2

DT Patent

LA German

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004043906	A2	20040527	WO 2003-EP12686	20031113 <--
	WO 2004043906	A3	20041007		
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BW, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, EG, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, SY, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW				
	RW: BW, GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	DE 10253331	A1	20040603	DE 2002-10253331	20021114 <--
	AU 2003283398	A1	20040603	AU 2003-283398	20031113 <--
	EP 1562893	A2	20050817	EP 2003-775352	20031113 <--
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				
	CN 1711234	A	20051221	CN 2003-80103209	20031113 <--
	JP 2006506479	T	20060223	JP 2004-551009	20031113 <--
	BR 2003016207	A	20060411	BR 2003-16207	20031113 <--
	US 2004241312	A1	20041202	US 2004-483668	20040727 <--
PRAI	DE 2002-10253331	A	20021114	<--	
	WO 2003-EP12686	W	20031113		

AB Use of 2E,4E-decadienoic acid isobutylamide (trans-pellitorin) (I) in the form of an aromatic substance, in particular a saliva stimulating aromatic substance for food, oral hygiene or gustatory prepsns. is claimed. Thus, a mixture of Et 2E,4Z-decadienoate, Chirazyme L-2, and isobutylamine was heated at 55° for 4 days to give 99.4% 2E,4Z-decadienoic acid isobutylamide, which was stirred 1 h with iodine in PhMe to give I in >95% purity. I food and oral hygiene compns. are given.

IC ICM C07C0231-00

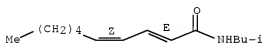
ICS A61K0007-16

CC 23-18 (Aliphatic Compounds)

Section cross-reference(s): 17, 62

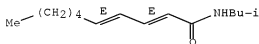
- ST dodecadienoate amidation isomerization; pellitorin prepn saliva  
stimulating substance food oral hygiene
- IT Chewing gum  
Dentifrices  
Flavoring materials  
Food additives  
Isomerization  
Mouthwashes  
Saliva  
(preparation and use of trans-pellitorin as an aromatic substance with  
salivation-stimulating activity)
- IT 639086-18-3P  
RL: BPN (Biosynthetic preparation); RCT (Reactant); BIOL (Biological  
study); PREP (Preparation); RACT (Reactant or reagent)  
(preparation and use of trans-pellitorin as an aromatic substance with  
salivation-stimulating activity)
- IT 18836-52-7F, trans-Pellitorin  
RL: COS (Cosmetic use); FFD (Food or feed use); SPN (Synthetic  
preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation and use of trans-pellitorin as an aromatic substance with  
salivation-stimulating activity)
- IT 639086-18-3P  
RL: BPN (Biosynthetic preparation); RCT (Reactant); BIOL (Biological  
study); PREP (Preparation); RACT (Reactant or reagent)  
(preparation and use of trans-pellitorin as an aromatic substance with  
salivation-stimulating activity)
- RN 639086-18-3 HCAPLUS  
CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2E,4Z)- (CA INDEX NAME)

Double bond geometry as shown.



- IT 18836-52-7F, trans-Pellitorin  
RL: COS (Cosmetic use); FFD (Food or feed use); SPN (Synthetic  
preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(preparation and use of trans-pellitorin as an aromatic substance with  
salivation-stimulating activity)
- RN 18836-52-7 HCAPLUS  
CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2E,4E)- (CA INDEX NAME)

Double bond geometry as shown.



- L111 ANSWER 7 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN  
AN 2004:101119 HCAPLUS [Full-text](#)  
DN 140:145102  
TI Flavorant aliphatic or aromatic unsatd. amide compounds for  
food use

IN Galopin, Christophe; Goeke, Andreas; Furrer,  
Stefan  
PA Givaudan SA, Switz.  
SO PCT Int. Appl., 15 pp.  
CODEN: PIXXD2  
DT Patent  
LA English  
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004011415	A1	20040205	WO 2003-CH500	20030723 <--
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NI, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
	RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
	AU 2003245800	A1	20040216	AU 2003-245800	20030723 <--
	EP 1525184	A1	20050427	EP 2003-737827	20030723 <--
	R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
	CN 1668578	A	20050914	CN 2003-816951	20030723 <--
	JP 2005533897	T	20051110	JP 2004-523719	20030723 <--
	IN 2004CN03060	A	20060217	IN 2004-CN03060	20041231 <--
	US 2005233042	A1	20051020	US 2005-522113	20050125 <--
PRAI	US 2002-398449P	P	20020725	<--	
	WO 2003-CH500	W	20030723	<--	
OS	MARPAT 140:145102				
AB	Use as a flavor ingredient of an aliphatic or aromatic unsatd. amide of formula (I, R''(CH)nCONHCH2C(R')(R''')(R''''')), where R' is H or OH, n is 1 or 2; R'' is RvCHCH(CH2)m when n is 2, m being 1,2 or 3; R''' and R'''' are H, C1-C4 alkyl, benzyl or form a 5- or 6-membered carbocyclic ring with the carbon to which they are attached; and Rv is alkyl or alkenyl is described. When n is 1, R'' may also be a Ph group.				
IC	ICM C07C0233-09				
	ICS C07C0233-20; A23L0001-226				
CC	17-6 (Food and Feed Chemistry)				
	Section cross-reference(s): 23				
ST	flavoring material aliph arom unsatd amide isobutyl undecatrienamide				
IT	Flavoring materials				
	Soups				
	(flavorant aliphatic or aromatic unsatd. amide compds. for food use)				
IT	Amides, biological studies				
	RL: FFD (Food or feed use); BIOL (Biological study); USES (Uses) (unsatd.); flavorant aliphatic or aromatic unsatd. amide compds. for food use)				
IT	652970-05-3P 652970-06-4P				
	RL: FFD (Food or feed use); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses) (flavorant aliphatic or aromatic unsatd. amide compds. for food use)				
IT	621-82-9, Cinnamic acid, reactions 4634-89-3, (Z)-Hex-4-enal 62872-62-2 67629-62-3				
	RL: RCT (Reactant); RACT (Reactant or reagent)				

(flavorant aliphatic or aromatic unsatd. amide compds. for food use)

IT 102-92-1P 6299-56-5P 652970-07-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(flavorant aliphatic or aromatic unsatd. amide compds. for food use)

IT 652970-06-3P 652970-06-4P

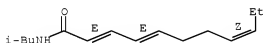
RL: EFD (Food or feed use); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)

(flavorant aliphatic or aromatic unsatd. amide compds. for food use)

RN 652970-06-3 HCAPLUS

CN 2,4,8-Undecatrienamide, N-(2-methylpropyl)-, (2E,4E,8Z)- (CA INDEX NAME)

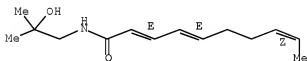
Double bond geometry as shown.



RN 652970-06-4 HCAPLUS

CN 2,4,8-Decatrienamide, N-(2-hydroxy-2-methylpropyl)-, (2E,4E,8Z)- (CA INDEX NAME)

Double bond geometry as shown.



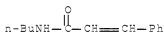
IT 6299-56-5P

RL: RCT (Reactant); SPN (Synthetic preparation); PREP (Preparation); RACT (Reactant or reagent)

(flavorant aliphatic or aromatic unsatd. amide compds. for food use)

RN 6299-56-5 HCAPLUS

CN 2-Propenamide, N-butyl-3-phenyl- (CA INDEX NAME)



# RETABLE

Referenced Author (RAU)	Year (RPY)	VOL (RVL)	PG (RPG)	Referenced Work (RWK)	Referenced File
Goeke, A	2002	1	1	JUS 2002081370 A1	1
Kikuzaki, H	1993	157	1329	BIOSCIENCE, BIOTECHN	HCAPLUS
Lion Corp	1985	1	1	JJP 60075424 A	HCAPLUS
Nakatani, N	1992	156	1759	BIOSCIENCE BIOTECHNO	HCAPLUS
Ramsewak, R	1999	151	1729	PHYTOCHEMISTRY	HCAPLUS

Sumitomo Chem Co Ltd	1979		JJP 54117476 A	HCAPLUS
Sumitomo Chem Co Ltd	1981		JJP 56087504 A	HCAPLUS

L111 ANSWER 8 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2004:2835 HCAPLUS Full-text

DN 140:59332

TI Amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant

IN Gatfield, Ian-Lucas; Ley, Jakob Peter; Foerstner, Jan; Krammer, Gerhard; Machinek, Arnold

PA Symrise G.m.b.H. &amp; Co. K.-G., Germany

SO PCT Int. Appl., 22 pp.

CODEN: PIXXD2

DT Patent

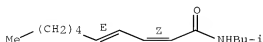
LA German

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2004000787	A2	20031231	WO 2003-EP6545	20030620 <--
	WO 2004000787	A3	20040805		
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MY, NI, NO, NZ, OM, PG, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
	RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
	DE 10227462	A1	20040108	DE 2002-10227462	20020620 <--
	DE 10227462	A9	20050728		
	AU 2003246570	A1	20040106	AU 2003-246570	20030620 <--
	EP 1517880	A2	20050330	EP 2003-760673	20030620 <--
	R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
	US 2005234132	A1	20051020	US 2004-518074	20041216 <--
PRAI	DE 2002-10227462	A	20020620	<--	
	WO 2003-EP6545	W	20030620	<--	
AB	A method for producing 2E,4Z-decadienoic acid isobutylamide (i.e., cis-pellitorin) is described as is its use as a pungent agent and a flavoring that generates heat, in foods, oral hygiene compns., or gourmet prepsns.				
IC	ICM C07C0231-00				
CC	23-18 (Aliphatic Compounds)				
	Section cross-reference(s): 17, 45, 62				
ST	decadienoic acid isobutylamide prepn flavoring agent odorant				
IT	Amides, preparation				
	RL: FFD (Food or feed use); PRP (Properties); SPN (Synthetic preparation); TEM (Technical or engineered material use); BIOL (Biological study); PREP (Preparation); USES (Uses)				
	(E,4Z-decadienoic acid isobutylamide; amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)				
IT	Flavoring materials				
	Food				
	Odor and Odorous substances				
	(amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)				
IT	Hygiene				
	(oral; amidation method for the production of cis-pellitorin and its use as				

- a flavoring agent and an odorant and in compns. for)
- IT Amidation  
(transamidation, enzymic; amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 73785-31-6 73785-32-7 175288-20-7 625092-39-9  
639086-19-4 639086-20-7 639086-21-8  
RL: FFD (Food or feed use); MOA (Modifier or additive use); BIOL (Biological study); USES (Uses)  
(additive; amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 18836-52-7P  
RL: FFD (Food or feed use); MOA (Modifier or additive use); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses)  
(amidation method for the production of cis-pellitorin and conversion to trans-pellitorin)
- IT 639086-18-3P  
RL: FFD (Food or feed use); PRP (Properties); RCT (Reactant); SPN (Synthetic preparation); TEM (Technical or engineered material use); BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent); USES (Uses)  
(amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 60-29-7, Diethylether, uses  
RL: NUU (Other use, unclassified); USES (Uses)  
(amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 1310-58-3, Potassium hydroxide, reactions  
RL: RCT (Reactant); RACT (Reactant or reagent)  
(amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 9001-62-1, Chirazyme L-2  
RL: CAT (Catalyst use); USES (Uses)  
(in an amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 3025-30-7  
RL: RCT (Reactant); RACT (Reactant or reagent)  
(in an amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- IT 175288-20-7 625092-39-9  
RL: FFD (Food or feed use); MOA (Modifier or additive use); BIOL (Biological study); USES (Uses)  
(additive; amidation method for the production of cis-pellitorin and its use as a flavoring agent and an odorant)
- RN 175288-20-7 HCAPLUS  
CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2Z,4E)- (CA INDEX NAME)

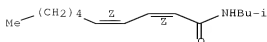
Double bond geometry as shown.



- RN 625092-39-9 HCAPLUS  
CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2Z,4Z)- (CA INDEX NAME)

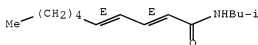
Double bond geometry as shown.





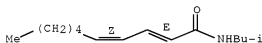
IT 19836-52-7P  
 RL: FFD (Food or feed use); MOA (Modifier or additive use); SPN  
 (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES  
 (Uses)  
 (amidation method for the production of cis-pellitorin and conversion to  
 trans-pellitorin)  
 RN 19836-52-7 HCAPLUS  
 CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2E,4E)- (CA INDEX NAME)

Double bond geometry as shown.



IT 639086-18-3P  
 RL: FFD (Food or feed use); PRP (Properties); RCT (Reactant);  
 SPN (Synthetic preparation); TEM (Technical or engineered material use);  
 BIOL (Biological study); PREP (Preparation); RACT (Reactant or reagent);  
 USES (Uses)  
 (amidation method for the production of cis-pellitorin and its use as a  
 flavoring agent and an odorant)  
 RN 639086-18-3 HCAPLUS  
 CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2E,4Z)- (CA INDEX NAME)

Double bond geometry as shown.

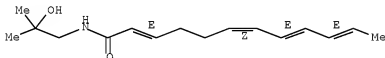


L111 ANSWER 9 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN  
 AN 2003:944238 HCAPLUS Full-text  
 DN 140:180285  
 TI Pungent and tingling compounds in Asian cuisine  
 AU Galopin, Christophe C.; Furrer, Stefan M.; Goëke,  
 Andreas  
 CS Givaudan Flavors R&D, Ingredient Systems, Cincinnati, OH, 45069,  
 USA  
 SO ACS Symposium Series (2004), 867(Challenges in Taste Chemistry and  
 Biology), 139-152  
 CODEN: ACSMC8; ISSN: 0097-6156  
 PB American Chemical Society  
 DT Journal; General Review  
 LA English  
 AB A review. Southern Asian cuisine is well known for its use of flavorful and  
 pungent spices. The sanshool chems., such as  $\alpha$ -hydroxy-sanshool from the

Japanese Sanshoo pepper and other Asian peppers, are particularly interesting because they not only give a hot sensation in the mouth cavity but also a tingling effect on the tongue. In order to understand the effect of the sanshool chems. the authors have synthesized a variety of derivs. Tasting of those derivs. provided information about Structure Activity Relationship (SAR) for the tingling effect exhibited by these chems. Based on this study the authors are able to propose a minimal structure required for the tingling effect. We also used this SAR knowledge to design stable compds. with potential tingling effect.

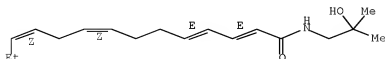
- CC 17-0 (Food and Feed Chemistry)  
 ST review Asian food additive Sanshoo bungeanool deriv pungency tingling; Sanshoo bungeanool deriv structure pungency tingling review  
 IT Taste  
     (pungency; pungent and tingling compds. in Asian cuisine)  
 IT Structure-activity relationship  
     (taste; pungent and tingling Sanshoo and bungeanool compds. in Asian cuisine)  
 IT Food functional properties  
     (tingling; pungent and tingling Sanshoo and bungeanool compds. in Asian cuisine)  
 IT 83883-10-7D,  $\alpha$ -Hydroxy-sanshool, derivs.  
     117568-40-8D, Bungeanool, derivs.  
     RL: BSU (Biological study, unclassified); FFD (Food or feed use); BIOL (Biological study); USES (Uses)  
     (pungent and tingling Sanshoo and bungeanool compds. in Asian cuisine)  
 IT 83883-10-7D,  $\alpha$ -Hydroxy-sanshool, derivs.  
     117568-40-8D, Bungeanool, derivs.  
     RL: BSU (Biological study, unclassified); FFD (Food or feed use); BIOL (Biological study); USES (Uses)  
     (pungent and tingling Sanshoo and bungeanool compds. in Asian cuisine)  
 RN 83883-10-7 HCAPLUS  
 CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-, (2E,6Z,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



- RN 117568-40-8 HCAPLUS  
 CN 2,4,8,11-Tetradecatetraenamide, N-(2-hydroxy-2-methylpropyl)-, (2E,4E,8Z,11Z)- (CA INDEX NAME)

Double bond geometry as shown.



# RETABLE

Referenced Author (RAU)	Year   VOL   PG   Referenced Work (RPY)   (RVL)   (RPG)   (RWK)   Referenced File
----------------------------	---

Bryant, B	1999  842  452	Brain Research	HCAPLUS
Chen, I	1999  52  357	Phytochemistry	HCAPLUS
Crombie, L	1952    4338	J Chem Soc	HCAPLUS
Crombie, L	1955    995	J Chem Soc	HCAPLUS
Crombie, L	1957    2760	J Chem Soc	HCAPLUS
Crombie, L	1985  26  2477	Tetrahedron Lett	HCAPLUS
Jacobson, M	1967  32  1646	J Org Chem	HCAPLUS
Mizutani, K	1988  36  2362	Chem Pharm Bull	HCAPLUS
Sonnet, P	1969  34  1147	J Org Chem	HCAPLUS
van der Linde, L	1985	EP 0173395 A1	HCAPLUS
Ward, J	1969  88  177	Recl Trav Chim Pays-	HCAPLUS
Xiong, Q	1997  46  1123	Phytochemistry	HCAPLUS

L111 ANSWER 10 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 2002:842774 HCAPLUS Full-text

DN 138:284551

TI Pungency and tingling: sensations and mechanisms of trigeminal chemical sensitivity

AU Bryant, Bruce; Mezzine, Igor

CS Monell Chemical Senses Center, Philadelphia, PA, 19104, USA

SO ACS Symposium Series (2002), 825(Chemistry of Taste), 202-212

CODEN: ACSMC8; ISSN: 0097-6156

PB American Chemical Society

DT Journal

LA English

AB Distinct from taste and olfaction, the trigeminal nerve is the third sensory pathway in the cranial sensory system that is sensitive to chemical stimuli. Trigeminal nerve endings in the nose and mouth contribute to flavor through the sensory modalities of touch, thermal sensation and pain. The best-characterized example of chemical induced trigeminal sensation is the pungency produced by hot peppers, the result of the activation of ion channels on pain-sensitive and thermally sensitive nerve fibers. Compds. commonly found in spices, food and beverages also elicit sensations other than pain. Menthol and other related compds. stimulate a subclass of thermal nerve endings to produce cooling. Yet other compds., stimuli as diverse as CO2 and fatty acids as well as some unsatd. alkylamides found in non-capsicum peppers and other plants, activate cooling-sensitive and tactile nerve endings. This particular combination of modalities gives rise to the novel tingling sensations associated with these stimuli.

CC 13-6 (Mammalian Biochemistry)

ST hydroxysanshool calcium pungency tingling trigeminal neurotransmission flavor

IT Taste

(pungency; pungency and tingling sensations and mechanisms of trigeminal chemical sensitivity)

IT 7440-70-2, Calcium, biological studies 23683-10-7,

Hydroxy- $\alpha$ -sanshool

RL: BSU (Biological study, unclassified); BIOL (Biological study)

(effects of hydroxy- $\alpha$ -sanshool on intraneuronal calcium and taste pungency in mechanisms of trigeminal chemical sensitivity)

IT 504-97-2,  $\alpha$ -Sanshool 7328-34-9 10076-00-3,  $\beta$ -Sanshool 18744-21-3 18836-52-7, Pellitorine 25394-57-4, Spilanthal 30361-33-2 65937-49-7 68125-01-9 73785-32-7 97465-69-5,

Hydroxy- $\beta$ -sanshool 252193-26-3, Hydroxy- $\epsilon$ -sanshool

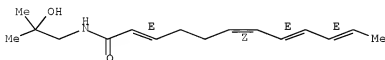
499136-10-6 499136-12-8

RL: BSU (Biological study, unclassified); BIOL (Biological study)

(pungency and tingling sensations and mechanisms of trigeminal chemical sensitivity)

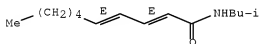
IT 83883-10-7, Hydroxy- $\alpha$ -sanshool  
 RL: BSU (Biological study, unclassified); BIOL (Biological study)  
 (effects of hydroxy- $\alpha$ -sanshool on intraneuronal calcium and  
 taste pungency in mechanisms of trigeminal chemical sensitivity)  
 RN 83883-10-7 HCAPLUS  
 CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-,  
 (2E,6Z,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



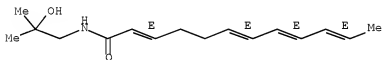
IT 18836-52-7, Pellitorine 97465-69-5, Hydroxy- $\beta$ -  
 sanshool 252193-26-3, Hydroxy- $\delta$ -sanshool  
 RL: BSU (Biological study, unclassified); BIOL (Biological study)  
 (pungency and tingling sensations and mechanisms of trigeminal chemical  
 sensitivity)  
 RN 18836-52-7 HCAPLUS  
 CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2E,4E)- (CA INDEX NAME)

Double bond geometry as shown.



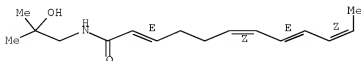
RN 97465-69-5 HCAPLUS  
 CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-,  
 (2E,6E,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



RN 252193-26-3 HCAPLUS  
 CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-,  
 (2E,6Z,8E,10Z)- (CA INDEX NAME)

Double bond geometry as shown.



## RETABLE

Referenced Author (RAU)	Year (RPY)	VOL (RVL)	PG (RPG)	Referenced Work (RWK)	Referenced File
-----	-----	-----	-----	-----	-----
Anon	1971		445	Fenaroli's Handbook	
Bryant, B	1999	842	452	Brain Res	HCAPLUS
Caterina, M	1997	389	816	Nature	HCAPLUS
Craig, A	1994	265	252	Science	MEDLINE
Duke, J	1985			CRC Handbook of Med	
Garnsworthy, R	1988	59	1116	J Neurophysiol	MEDLINE
Green, B	1992	17	435	Chemical Senses	HCAPLUS
Greger, H	1984	50	366	Planta Medica	HCAPLUS
Hegnauer, R	1977			The Biology and Chem	
Holzer, P	1991	43	143	Pharmacol Rev	HCAPLUS
Jacobson, M	1948	70	4234	J Am Chem Soc	HCAPLUS
Kashiwada, Y	1997	44	1125	Phytochem	HCAPLUS
Liu, L	1996	76	1858	J Neurophysiol	HCAPLUS
Martenson, M	1997	761	71	Brain Res	HCAPLUS
Schmeiz, M	1997	17	8003	J Neurosci	HCAPLUS
Walpole, C	1993	36	2381	J Med Chem	HCAPLUS
Yasuda, I	1981	29	1791	Chem Pharm Bull	HCAPLUS

L111 ANSWER 11 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 1999:297270 HCAPLUS [Full-text](#)

DN 130:329049

TI Trigeminal sensory stimuli and animal repellents from plants

IN Bryant, Bruce P.; Mezine, Igor A.; Epple, Gisela M.

PA Monell Chemical Senses Center, USA

SO PCT Int. Appl., 28 pp.

CODEN: PIXXD2

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
	-----	----	-----	-----	-----
PI	WO 9921425	A1	19990506	WO 1998-US22537	19981023 <--
	W: CA, JP				
	RW: AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL,				
	PT, SE				

PRAI US 1997-957112 A 19971024 &lt;--

AB Novel uses for compds. isolated from the fruit of Xanthoxylum and echinacea species, and similar compds. from other spice and flowering species, and the oil exts. from which they are isolated, are disclosed. The novel uses include flavor enhancers, additives for oral, hair, and skin care products, and animal repellents. An Et acetate extract of Xanthoxylum fruit was prepared, then evaporated to obtain an oil-like black-brown liquid of which hydroxy- $\alpha$ -sanshool (I) was separated. I induced increase in calcium in neurons that were sensitive or insensitive to capsaicin. Twenty hour food-deprived rats consumed significantly less rat chow that had been treated with a crude vegetable oil suspension of Xanthoxylum extract than rats chow treated with a similar oil suspension of an extract made of equal weight of cinnamon.

IC ICM A01N0065-00

CC 62-5 (Essential Oils and Cosmetics)

Section cross-reference(s): 1, 5, 17

IT Bird (Aves)

(food repellents for; trigeminal sensory stimuli and animal repellents from plants)

IT 82883-10-7, Hydroxy- $\alpha$ -sanshool

RL: BAC (Biological activity or effector, except adverse); BOC (Biological

occurrence); BSU (Biological study, unclassified); BUU (Biological use, unclassified); FFD (Food or feed use); BIOL (Biological study); OCCU (Occurrence); USES (Uses)

(trigeminal sensory stimuli and animal repellents from plants)

IT 83883-10-7, Hydroxy- $\alpha$ -sanshool

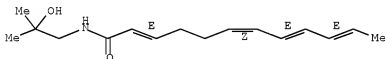
RL: BAC (Biological activity or effector, except adverse); BOC (Biological occurrence); BSU (Biological study, unclassified); BUU (Biological use, unclassified); FFD (Food or feed use); BIOL (Biological study); OCCU (Occurrence); USES (Uses)

(trigeminal sensory stimuli and animal repellents from plants)

RN 83883-10-7 HCAPLUS

CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-, (2E,6Z,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



# RETABLE

Referenced Author (RAU)	Year (RPY)	VOL (RVL)	PG (RPG)	Referenced Work (RWK)	Referenced File
Wunderlich	1995			JUS 5401502 A	HCAPLUS

L111 ANSWER 12 OF 13 HCAPLUS COPYRIGHT 2008 ACS on SIN

AN 1997:155786 HCAPLUS [Full-text](#)

DN 126:261550

TI Amides of the fruit of Zanthoxylum spp

AU Kashiwada, Yoshiki; Ito, Chikashi; Katagiri, Hitoshi; Mase, Izumi; Komatsu, Katsuko; Namba, Tsuneo; Ikeshiro, Yasumasa

CS Niigata College Pharmacy, Niigata, 950-21, Japan

SO Phytochemistry (1997), 44(6), 1125-1127

CODEN: PYTCAS; ISSN: 0031-9422

PB Elsevier

DT Journal

LA English

AB Examination of the amide constituents in Budo-Zanthoxylum fruit, the most traded com. Zanthoxylum fruit in the Japanese market, has led to the isolation of a new amide, along with  $\alpha$ -,  $\beta$ -,  $\gamma$ -, hydroxy- $\alpha$ -, hydroxy- $\beta$ - and hydroxy- $\gamma$ -sanshools. The structure of the new amide was assigned as (2E,4E,8E,10E,12E)-N-isobutyl-2,4,8,10,12-tetradecapentaenamide by spectral examination

CC 11-1 (Plant Biochemistry)

Section cross-reference(s): 17, 26

IT 504-97-2P,  $\alpha$ -Sanshool 10076-00-3P,  $\beta$ -Sanshool 78886-65-4P,  $\gamma$ -Sanshool 78886-66-5P, Hydroxy- $\gamma$ -sanshool

83883-10-7P, Hydroxy- $\alpha$ -sanshool 97465-69-5P

RL: BOC (Biological occurrence); BSU (Biological study, unclassified); PUR (Purification or recovery); BIOL (Biological study); OCCU (Occurrence); PREP (Preparation)

(amides of Zanthoxylum fruit)

IT 78886-66-5P, Hydroxy- $\gamma$ -sanshool 83883-10-7P,

Hydroxy- $\alpha$ -sanshool 97465-69-5P

RL: BOC (Biological occurrence); BSU (Biological study, unclassified); PUR (Purification or recovery); BIOL (Biological study); OCCU (Occurrence);

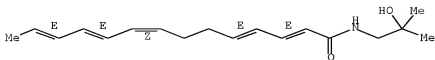
## PREP (Preparation)

(amides of Zanthoxylum fruit)

RN 78886-66-5 HCAPLUS

CN 2,4,8,10,12-Tetradecapentaenamide, N-(2-hydroxy-2-methylpropyl)-,  
(2E,4E,8Z,10E,12E)- (CA INDEX NAME)

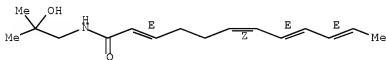
Double bond geometry as shown.



RN 83883-10-7 HCAPLUS

CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-,  
(2E,6Z,8E,10E)- (CA INDEX NAME)

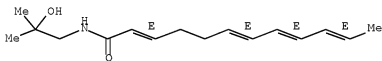
Double bond geometry as shown.



RN 97465-69-5 HCAPLUS

CN 2,6,8,10-Dodecatetraenamide, N-(2-hydroxy-2-methylpropyl)-,  
(2E,6E,8E,10E)- (CA INDEX NAME)

Double bond geometry as shown.



L111 ANSWER 13 OF 13 HCAPLUS COPYRIGHT 2008 ACS on STN

AN 1993:190280 HCAPLUS [Full-text](#)

DN 118:190280

TI Amides from supercritical fluid extracts of muntok pepper

AU Kollmannsberger, H.; Nitz, S.

CS Inst. Lebensmitteltechnol. Anal. Chem., Tech. Univ. Muenchen, Freising,  
W-8050, GermanySO Chemie, Mikrobiologie, Technologie der Lebensmittel (1992),  
14(3/4), 87-94

CODEN: CMTLBX; ISSN: 0366-7154

DT Journal

LA German

AB In supercrit. fluid exts. of Muntok pepper, 21 piperidides, 7 pyrrolidides,  
and 7 isobutylamides of various saturated and unsatd. fatty acids and 3,4-  
methylenedioxybenzyl-substituted carbonic acids were separated by gas  
chromatog. (GC) and identified by mass spectrometry (GC-MS) and in some cases

by IR spectroscopy (GC-FTIR). Their sensorial contributions and physiol. effects are briefly discussed.

CC 17-6 (Food and Feed Chemistry)

IT Amides, biological studies

RL: BIOL (Biological study)

(of Muntok pepper flavor)

IT Flavor

(of Muntok pepper, piperidides and pyrrolidides and isobutylamides of)

IT 94-62-2, Piperin 94-62-2 583-34-6, Piperettin 618-42-8,  
Acetylpiiperidide 2591-86-8, Formylpiiperidide 4629-02-1,  
Hexadecanoylpiiperidide 5299-66-1, Dodecanoylpiiperidide  
18836-52-7 23512-46-1 24738-51-0 25924-78-1, Piperylin  
27845-72-3 30505-89-6 30505-92-1 42997-42-2 54794-69-3  
54794-70-6 56630-42-3, 9-Octadecanoylpiiperidide 65937-45-3  
78910-33-5 82857-82-7 91487-76-2 117137-69-6 145398-89-6  
145398-91-0 145398-95-4 145427-76-5 147030-02-2 147030-03-3  
147030-04-4 147030-05-5 147030-06-6 147030-09-9 147030-13-5

RL: BIOL (Biological study)

(of Muntok pepper aroma)

IT 18836-52-7

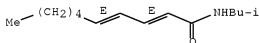
RL: BIOL (Biological study)

(of Muntok pepper aroma)

RN 18836-52-7 HCAPLUS

CN 2,4-Decadienamide, N-(2-methylpropyl)-, (2E,4E)- (CA INDEX NAME)

Double bond geometry as shown.



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(FILE 'HOME' ENTERED AT 13:33:28 ON 26 MAR 2008)

SET COST OFF

FILE 'HCAPLUS' ENTERED AT 13:33:56 ON 26 MAR 2008

L1 1 S US20050233042/PN OR (US2005-522113# OR WO2003-CH500)/AP,PRN  
E GALOPIN/AU  
L2 30 S E4,E5,E8-E10  
E GOEKE/AU  
L3 28 S E3,E6  
E GEOKE/AU  
E GOKE/AU  
E FURRER/AU  
E FURRER S/AU  
L4 30 S E3-E6  
E FUERRER/AU  
E GIVAUDAN/CO  
L5 1348 S E3-E96  
E E82+ALL  
L6 1845 S E2+RT OR E2-E50/PA,CS  
SEL RN L1

FILE 'REGISTRY' ENTERED AT 13:36:38 ON 26 MAR 2008

L7 9 S E1-E9



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L8      3 S L7 AND N/ELS AND C>=10 NOT P/ELS

FILE 'HCAPLUS' ENTERED AT 13:39:06 ON 26 MAR 2008
L9      68 S L2-L4 NOT L1

FILE 'REGISTRY' ENTERED AT 13:39:19 ON 26 MAR 2008

FILE 'HCAPLUS' ENTERED AT 13:39:19 ON 26 MAR 2008
L10     TRA L9 1- RN :      837 TERMS

FILE 'REGISTRY' ENTERED AT 13:39:21 ON 26 MAR 2008
L11     837 SEA L10
L12     16 S L11 AND 1/O AND 1/N AND 4/ELC.SUB
L13     24 S L11 AND 2/O AND 1/N AND 4/ELC.SUB
L14     2 S L13 AND (C18H29NO2 OR C16H25NO2) NOT C6/ES
L15     5 S L8,L14
        E C1317 N O/MF
        E C13H17NO/MF
L16     2287 S E3 AND 46.150.18/RID
L17     51 S L16 AND 2 PROPENAMIDE
L18     19 S L17 AND 3 PHENYL
L19     6 S L18 AND METHYLPROPYL
L20     3 S L19 AND 2 METHYLPROPYL
        E C11H19NO/MF
L21     1899 S E3
L22     1681 S L21 AND NR>=1
L23     218 S L21 NOT L22
L24     4 S L23 AND METHYLPROPYL
L25     1 S L24 AND HEPTADIENAMIDE
        E C10H17NO2/MF
L26     2613 S E3
L27     2201 S L26 AND NR>=1
L28     412 S L26 NOT L27
L29     193 S L28 NOT ESTER
L30     177 S L29 NOT CYAN?
L31     148 S L30 NOT NITRILE
L32     132 S L31 NOT NITRO
L33     115 S L32 NOT METHOXY
        E C14H25NO/MF
L34     926 S E3
L35     819 S L34 AND NR>=1
L36     107 S L34 NOT L35
L37     97 S L36 NOT (NITRO OR NITRILE OR CYAN? OR ESTER)
L38     13 S L37 AND METHYLPROPYL
L39     6 S L38 AND 2 4
        E C14H25NO2/MF
L40     1136 S E3
L41     952 S L40 AND NR>=1
L42     184 S L40 NOT L41
L43     62 S L42 NOT (NITRO OR NITRILE OR CYAN? OR ESTER)
L44     60 S L43 NOT ACETATE
L45     52 S L44 NOT OXIME
        E C10H17NO2/MF
L46     15 S L15,L20,L25,L39
        SEL RN 11 14
L47     13 S L46 NOT E1-E2
        E C14H23NO2/MF
L48     1926 S E3
L49     1833 S L48 AND NR>=1
L50     93 S L48 NOT L49

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L51      92 S L50 NOT HYDROXYETHYL
L52      28 S L51 NOT (NITRO OR NITRILE OR CYAN? OR ESTER)
L53      26 S L52 NOT ?NITRIL?/CNS
L54      1 S L53 AND DECATRIENAMIDE AND HYDROXY AND METHYLPROPYL
L55      1 S L50 AND METHYLPROPYL AND HYDROXY
L56      13 S L47,L54,L55
L57      2 S L46 NOT L56
L58      STR
L59      50 S L58
L60      42913 S L58 FUL
L61      STR L58
L62      50 S L61 CSS SAM SUB=L60
L63      1639 S L61 CSS FUL SUB=L60
          SAV TEMP L63 DEES522A/A
L64      STR L58
L65      4 S L64 CSS SAM SUB=L63
L66      94 S L64 CSS FUL SUB=L63
          SAV TEMP L66 DEES522B/A
L67      3 S L66 AND NC>=2
L68      2 S L67 NOT C42H70O35
L69      1 S L67 NOT L68
L70      93 S L66 NOT L69
L71      103 S L56,L70
L72      STR L64
L73      2 S L72 SAM SUB=L63
L74      22 S L72 FUL SUB=L63
          SAV TEMP L74 DEES522C/A
L75      4 S L74 NOT L71
L76      107 S L71,L74
          SAV TEMP L76 DEES522D/A

FILE 'HCAPLUS' ENTERED AT 14:23:18 ON 26 MAR 2008
L77      434 S L76
L78      2 S L77 AND L1-L6
L79      2 S L77 AND GIVAUDAN?/CO,PA,CS
L80      2 S L78,L79
          SEL RN

FILE 'REGISTRY' ENTERED AT 14:24:01 ON 26 MAR 2008
L81      11 S E1-E11
L82      2 S L81 AND N/ELS NOT L76
L83      1 S L82 AND C13H17NO

FILE 'HCAPLUS' ENTERED AT 14:25:09 ON 26 MAR 2008
L84      19 S L83
L85      1 S L1-L6 AND L84
L86      2 S L80,L85
L87      446 S L77,L84
L88      309 S L87 AND PY<=2003 NOT P/DT
L89      42 S L87 AND (PD<=20030723 OR PRD<=20030723) AND P
L90      351 S L88,L89
L91      7 S L90 AND (L77 OR L84) (L)FFD/RL
          E FLAVOR/CT
L92      31940 S E3-E8 OR E22-E35
          E E3+ALL
L93      29719 S E2,E3,E5-E9
          E E11+ALL
L94      21825 S E2,E9,E10
          E E12+ALL
L95      6113 S E4,E5

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L96      7 S L90 AND L92-L95
L97      9 S L86,L91,L96
          E TASTE/CT
L98      8336 S E3-E15
          E E3+ALL
L99      7821 S E4
          E E7+ALL
L100     1629 S E15+OLD
L101      3 S L90 AND L98-L100
L102     11 S L97,L101
L103     11 S L102 AND (TASTE OR FLAVOR? OR FLAVOUR?)
L104     11 S L102 AND (FEED? OR FOOD?)/CW,CT,SC, SX,BI
L105     11 S L102-L104
L106     18 S L90 AND (FEED? OR FOOD?)/CW,CT,SC, SX,BI
L107     12 S L90 AND (TASTE OR FLAVOR? OR FLAVOUR?)
L108      9 S L106,L107 NOT L105
          SEL AN 2 4
L109      2 S L108 AND E1-E4
L110     13 S L105,L109
L111     13 S L110 AND L1-L6,L77-L80,L84-L110

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FILE 'REGISTRY' ENTERED AT 14:36:46 ON 26 MAR 2008

FILE 'HCAPLUS' ENTERED AT 14:37:37 ON 26 MAR 2008

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